

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090921\
 Data File : BN016279.D
 Acq On : 10 Sep 2021 05:44
 Operator : CG/JU
 Sample : M3620-03
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BP-TT-VPB-RW7-GW-640-642

Manual Integrations
 APPROVED

mohammad
 9/10/2021 3:24:04 PM

Quant Time: Sep 10 06:12:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN090921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 09 15:19:08 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	12558	0.400	ng	0.00
7) Naphthalene-d8	10.660	136	64556	0.400	ng	0.00
13) Acenaphthene-d10	14.484	164	36991	0.400	ng	-0.01
19) Phenanthrene-d10	17.237	188	73558	0.400	ng	0.00
29) Chrysene-d12	21.429	240	69928	0.400	ng	0.00
35) Perylene-d12	23.810	264	47932	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.457	112	2456	0.077	ng	0.00
5) Phenol-d6	7.034	99	2027	0.051	ng	0.00
8) Nitrobenzene-d5	9.013	82	13940	0.269	ng	0.00
11) 2-Methylnaphthalene-d10	12.243	152	26670	0.265	ng	0.00
14) 2,4,6-Tribromophenol	15.982	330	2871	0.189	ng	0.00
15) 2-Fluorobiphenyl	13.114	172	38793	0.268	ng	0.00
27) Fluoranthene-d10	19.271	212	74226	0.343	ng	0.00
31) Terphenyl-d14	19.867	244	70770	0.403	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.373	88	9781m	2.195	ng	
9) Naphthalene	10.698	128	9373	0.054	ng	97
12) 2-Methylnaphthalene	12.318	142	9558	0.082	ng	99
17) Acenaphthene	14.548	154	5149	0.048	ng	# 73
18) Fluorene	15.537	166	13336	0.100	ng	# 76
24) Pentachlorophenol	16.896	266	833	0.235	ng	97
25) Phenanthrene	17.273	178	82103	0.384	ng	# 92
28) Fluoranthene	19.296	202	6500	0.026	ng	# 91
30) Pyrene	19.664	202	12784	0.054	ng	# 72
34) Bis(2-ethylhexyl)phtha...	21.334	149	38748	0.289	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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